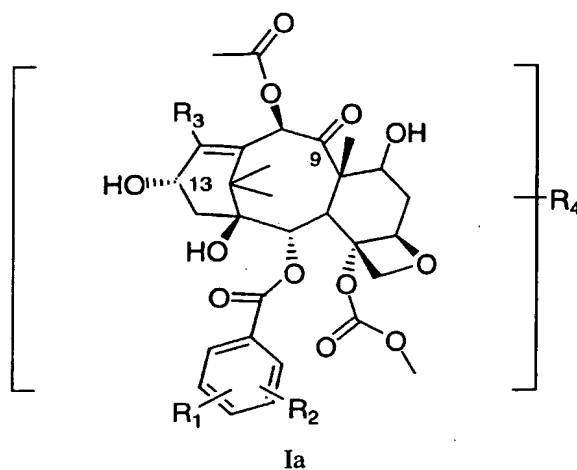


5 What is claimed is:

1. A metabolite of 3'-*tert*-butyl-3'-*N-tert*-butyloxycarbonyl-4-deacetyl-3'-dephenyl-3'-*N*-debenzoyl-4-O-methoxycarbonyl-paclitaxel of formula Ia or a pharmaceutically acceptable salt, solvate, hydrate or prodrug thereof

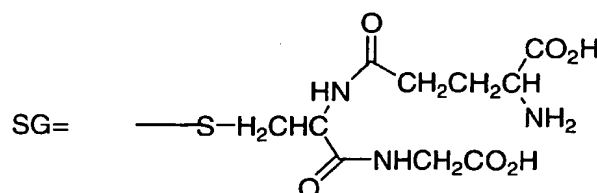
10



<u>Metabolite</u> <u>Code</u>	<u>R1</u>	<u>R2</u>	<u>R3</u>	<u>R4</u>	<u>C9</u>
M1	SG	H	CH <sub>3</sub>	OH	C=O
M2	SG	OH	CH <sub>3</sub>	H	C=O
M4 & M5	SG	H	CH <sub>3</sub>	H	C=O
M6	OCH <sub>3</sub>	OH	CH <sub>3</sub>	OH	C=O
M7	OH	OCH <sub>3</sub>	CH <sub>3</sub>	OH	C=O
M8	OH	H	CH <sub>3</sub>	OH	C=O
M8A	H	H	CH <sub>3</sub>	(OH) <sub>2</sub>	C=O
M9	H	H	CH <sub>3</sub>	H	(CH)OH

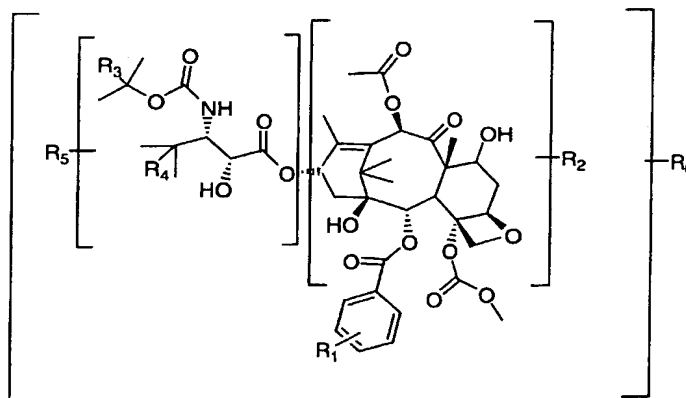
M10	H	H	CH <sub>3</sub>	OH	C=O
M11	H	H	COOH	H	C=O
M12	H	H	CH <sub>3</sub>	OH	C=O
M13	H	H	CH <sub>3</sub>	H	C=O

and



5

2. A metabolite of 3'-*tert*-butyl-3'-*N-tert*-butyloxycarbonyl-4-deacetyl-3'-dephenyl-3'-*N*-debenzoyl-4-*O*-methoxycarbonyl-paclitaxel of formula Ib or a pharmaceutically acceptable salt, solvate or prodrug thereof



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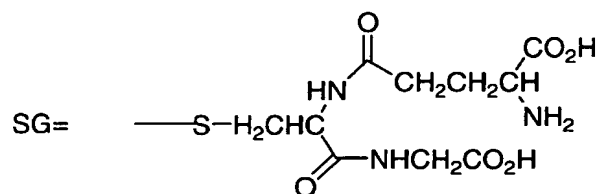
Ib

wherein the substituents are as defined in the following table

<u>Metabolite Code</u>	<u>R<sub>1</sub></u>	<u>R<sub>2</sub></u>	<u>R<sub>3</sub></u>	<u>R<sub>4</sub></u>	<u>R<sub>5</sub></u>	<u>R<sub>6</sub></u>
M14	H	H	CH <sub>3</sub>	CH <sub>3</sub>	H	2O
M15	SG	H	CH <sub>3</sub>	CH <sub>3</sub>	H	H
M16	H	H	CH <sub>3</sub>	CH <sub>3</sub>	H	2O

M17	H	H	CH <sub>3</sub>	CH <sub>3</sub>	OH	H
M18	H	OH	CH <sub>3</sub>	CH <sub>3</sub>	OH	H
M19	H	H	CH <sub>3</sub>	CH <sub>3</sub>	OH	H
M20	H	H				H
M21	H	H				H
M22	H	H	CH <sub>3</sub>	COOH	H	H
M23	H	H	CH <sub>3</sub>	CH <sub>3</sub>	OH	H
M24	H	H	COOH	CH <sub>3</sub>	H	H

and

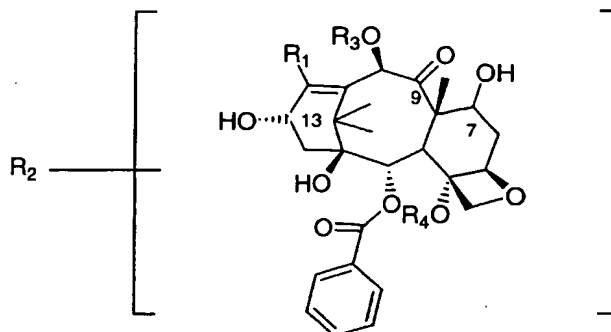


wherein the side chains on M20 and M21 are as shown below:

5

M20	
M21	

3. A metabolite of 3'-*tert*-butyl-3'-*N-tert*-butyloxycarbonyl-4-deacetyl -3'-dephenyl-3'-*N*-debenzoyl-4-*O*-methoxycarbonyl-paclitaxel of formula Ic or a  
10 pharmaceutically acceptable salt, solvate or prodrug thereof

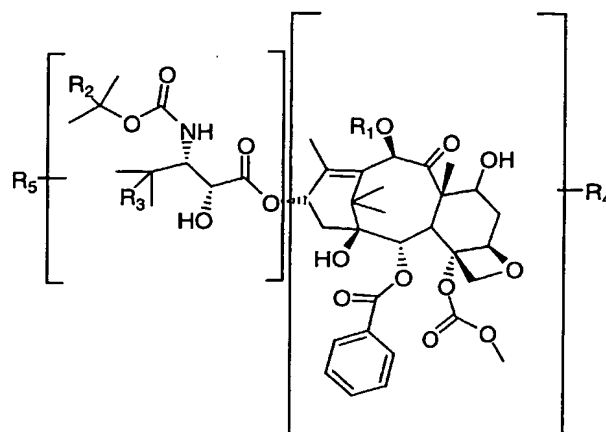


1c

5 wherein the substituents are as defined in the following table:

<u>Metabolite Code</u>	<u>R<sub>1</sub></u>	<u>R<sub>2</sub></u>	<u>R<sub>3</sub></u>	<u>R<sub>4</sub></u>	<u>C9</u>	<u>C13</u>
M9	CH <sub>3</sub>	H	CO(CH <sub>3</sub> )	CO(OCH <sub>3</sub> )	(CH)OH	(CH)OH
M10	CH <sub>3</sub>	OH	CO(CH <sub>3</sub> )	CO(OCH <sub>3</sub> )	C=O	(CH)OH
M10A	CH <sub>3</sub>	H	H	CO(OCH <sub>3</sub> )	C=O	(CH)OH
M11	COOH	H	CO(CH <sub>3</sub> )	CO(OCH <sub>3</sub> )	C=O	(CH)OH
M10B	CH <sub>3</sub>	OH	CO(CH <sub>3</sub> )	CO(OCH <sub>3</sub> )	C=O	(CH)OH
M10C	CH <sub>3</sub>	H	CO(CH <sub>2</sub> OH)	CO(OCH <sub>3</sub> )	C=O	(CH)OH
M12	CH <sub>3</sub>	OH	CO(CH <sub>3</sub> )	CO(OCH <sub>3</sub> )	C=O	(CH)OH
M13	CH <sub>3</sub>	H	CO(CH <sub>3</sub> )	CO(OCH <sub>3</sub> )	C=O	(CH)OH
M13A	CH <sub>3</sub>	H	H	CO(OCH <sub>3</sub> )	C=O	C=O
M13B	CH <sub>3</sub>	OH	CO(CH <sub>3</sub> )	H	C=O	(CH)OH
M13C	CH <sub>3</sub>	H	CO(CH <sub>3</sub> )	CO(OCH <sub>3</sub> )	C=O	(CH)OH
M13D	CH <sub>3</sub>	H	CO(CH <sub>3</sub> )	CO(OCH <sub>3</sub> )	C=O	C=O

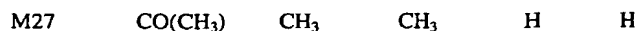
4. A metabolite of 3'-*tert*-butyl-3'-*N-tert*-butoxycarbonyl-4-deacetyl-3'-dephenyl-3'-*N*-debenzoyl-4-*O*-methoxycarbonyl-paclitaxel of formula Id or a pharmaceutically acceptable salt, solvate or prodrug thereof.



1d

wherein the substituents are as defined in the following table:

<u>Metabolite Code</u>	<u>R<sub>1</sub></u>	<u>R<sub>2</sub></u>	<u>R<sub>3</sub></u>	<u>R<sub>4</sub></u>	<u>R<sub>5</sub></u>
M15B	CO(CH <sub>3</sub> )	CH <sub>2</sub> OH or COOH	COOH or CH <sub>2</sub> OH	H	H
M17	CO(CH <sub>3</sub> )	CH <sub>3</sub>	CH <sub>3</sub>	H	OH
M18B	H	CH <sub>3</sub>	COOH	H	H
M19	CO(CH <sub>3</sub> )	CH <sub>3</sub>	CH <sub>3</sub>	H	OH
M19A	H	COOH	CH <sub>3</sub>	H	H
M22	CO(CH <sub>3</sub> )	CH <sub>3</sub>	COOH	H	H
M23	CO(CH <sub>3</sub> )	CH <sub>3</sub>	CH <sub>3</sub>	H	OH
M24	CO(CH <sub>3</sub> )	COOH or CH <sub>3</sub>	CH <sub>3</sub> or COOH	H	H
M23A	CO(CH <sub>3</sub> )	CH <sub>3</sub>	CH <sub>3</sub>	OH or H	H or OH
M23B	CO(CH <sub>3</sub> )	CH <sub>3</sub>	CH <sub>3</sub>	OH	H
M23C	CO(CH <sub>3</sub> )	CH <sub>3</sub>	CH <sub>3</sub>	OH or H	H or OH
M26	H	CH <sub>3</sub>	CH <sub>3</sub>	H	H
M23D	CO(CH <sub>3</sub> )	CH <sub>3</sub>	CH <sub>3</sub>	OH	H



5. A pharmaceutical composition comprising a metabolite according to Claim 1 or a pharmaceutically acceptable salt, solvate or prodrug thereof, and a pharmaceutically acceptable carrier, vehicle or diluent.
- 5 6. A pharmaceutical composition comprising a metabolite according to Claim 2 or a pharmaceutically acceptable salt, solvate or prodrug thereof, and a pharmaceutically acceptable carrier, vehicle or diluent.
- 10 7. A pharmaceutical composition comprising a metabolite according to Claim 3 or a pharmaceutically acceptable salt, solvate or prodrug thereof, and a pharmaceutically acceptable carrier, vehicle or diluent.
8. A pharmaceutical composition comprising a metabolite according to Claim 4 or a pharmaceutically acceptable salt, solvate or prodrug thereof, and a pharmaceutically acceptable carrier, vehicle or diluent.
- 15 9. A method for inhibiting tumor growth in a mammalian host which comprises administering to said mammal a tumor-growth inhibiting amount of a compound as defined in Claim 1.
- 20 10. A method for inhibiting tumor growth in a mammalian host which comprises administering to said mammal a tumor-growth inhibiting amount of a compound as defined in Claim 2.
- 25 11. A method for inhibiting tumor growth in a mammalian host which comprises administering to said mammal a tumor-growth inhibiting amount of a compound as defined in Claim 3.

12. A method for inhibiting tumor growth in a mammalian host which comprises administering to said mammal a tumor-growth inhibiting amount of a compound as defined in Claim 4.